CLAIMS

- 5 1. A compound comprising at least one carbon bearing:
 - an amine function;
 - an allyl or propargyl radical;
 - a difluoromethylene group;
 and
- 10 a hydrogen or a hydrocarbon-based radical advantageously chosen from those which are electron-donating or weakly electron-withdrawing ($\sigma_p \leq 0.2$, advantageously to 0.1) radicals.
- 15 2. A compound as claimed in claim 1, characterized in that the number of carbons of said compound is at most 30 carbon atoms, advantageously at most 20 carbon atoms.
- 20 3. A method for synthesizing a compound of formula I:

where:

- Rf represents a carbon radical bearing a difluoromethylene group providing the link with the rest of the molecule, advantageously of at most 15, preferably of at most 10, carbon atoms;
- R₁ represents a hydrogen, an alkyl, including aralkyl, radical, preferably of 1 or 2 carbon atoms, or one of the specific radicals subsequently specified;
- R₂ represents a hydrogen, an alkyl, including aralkyl, radical, preferably of 1 or 2 carbon atoms, or an aryl radical;

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- R₃ represents a hydrogen or a hydrocarbon-based radical, such as alkyl, including aralkyl, preferably of 1 or 2 carbon atoms, or an aryl radical, or forms, with R₄, an additional double bond so as to convert the allyl radical into a propargyl radical;
- R₄ represents a hydrogen or a hydrocarbon-based radical such as aryl or alkyl, including aralkyl, preferably of 1 or 2 carbon atoms, or, with R₃, forms an additional double bond so that the ethylenic bond becomes acetylenic, making it possible to go from an allyl radical to a propargyl radical;
- R₅ represents a hydrogen or a hydrocarbon-based radical such as an aryl radical or an alkyl, including aralkyl, radical, preferably of 1 or 2 carbon atoms;
 - R_5 and R_4 may be fractions of said "Ar" group above, such that R_5 and R_4 , and also the carbon which bears them, form a radical Ar;
 - it being possible for one of R_1 , R_2 , R_4 , R_3 and R_5 to be, in addition, chosen from specific trivalent, nitrile or acid functions, optionally and preferably in esterified form;
- R' may be a hydrogen, a protective group, an aryl or an alkyl, including aralkyl, advantageously a chiral alkyl or aralkyl;
 - R" is an allyl radical, a hydrogen or a metal cation, or a fraction of metal cations, when the metal is polyvalent;

by means of the action of an allyl organometallic on an imine bearing difluoromethylene groups.

- 4. The method as claimed in claim 3, characterized in that said organometallic is prepared in situ, according to a "Barbier" technique.
 - 5. The method as claimed in claims 3 and 4, characterized in that the imine is of formula

(II):

- where R is chosen from the halogens and from the hydrocarbon-based radicals chosen from those which are electron-donating or weakly electron-withdrawing radicals, and Rf has the same value as above and where R' is advantageously an alkyl, including aralkyl, including chiral, protective group.
- 6. The method as claimed in claims 3 to 5, characterized in that the allyl radical of said organometallic corresponds to formula III:

- 7. The method as claimed in claims 3 to 6, characterized in that the reaction is carried out in a polar aprotic solvent for which the donor number is at least equal to 10, advantageously at least equal to 20.
- 25 8. The method as claimed in claims 3 to 7, characterized in that it also comprises a step of N-allylation, by means of the action of an allyl derivative of formula IV on the free amine.
- 30 9. The method as claimed in claims 3 to 8, characterized in that the Rf group corresponds to the formula below:

- 33 -

$GEA-(CX_2)_p-$

where:

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- the X, which may be similar or different, represent a chlorine, a fluorine or a radical of formula C_nF_{2n+1} , with n an integer at most equal to 5, preferably to 2, with the condition that at least one of the X is fluorine, which fluorine is advantageously borne by the carbon bearing the open bond;
- p represents an integer at most equal to 2, i.e. 1 or 2;
 - GEA represents an electron-withdrawing group (i.e. sigma p greater than zero, advantageously than 0.1, preferably than 0.2) the possible functions of which are inert under the reaction conditions, advantageously fluorine or a perfluorinated residue of formula C_nF_{2n+1} , with n an integer at most equal to 8, advantageously to 5;
- the total number of carbons of Rf being advantageously between 1 and 15, preferably between 1 and 10.
- 10. The method as claimed in claim 9, characterized in that the total number of carbons of the radical Rf is between 1 and 14, advantageously between 1 and 10, more preferably from 1 to 4.
- The 11. method as claimed in claims 9 and 10, 30 characterized in that the radical Rf is a radical sensu stricto, i.e. it corresponds to the formula C_{ν} , $F_{\nu+1}$, where ν is an integer ranging from 1 to 10, advantageously of at most 4, most commonly of at most 2.
 - 12. The method as claimed in claims 9 to 11, characterized in that the radical Rf is chosen from the difluoromethyl (CHF_2) radical and the trifluoromethyl radical, the latter being

preferred.

13. The use, as substrate for cyclizing metathesis, of the compounds of formula (I):

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where:

- Rf represents a carbon radical bearing a difluoromethyl group providing the link with the rest of the molecule, advantageously of at most 15, preferably of at most 10, carbon atoms;
- R₁ represents a hydrogen, an alkyl, including aralkyl, radical, preferably of 1 or 2 carbon atoms, or one of the specific radicals subsequently specified;
 - R_2 represents a hydrogen, an alkyl, including aralkyl, radical, preferably of 1 or 2 carbon atoms, or an aryl radical;
 - R₃ represents a hydrogen or a hydrocarbon-based radical, such as alkyl, including aralkyl, preferably of 1 or 2 carbon atoms, or an aryl radical, or forms, with R₄, an additional double bond so as to convert the allyl radical into a propargyl radical;
 - R₄ represents a hydrogen or a hydrocarbon-based radical such as aryl or alkyl, including aralkyl, preferably of 1 or 2 carbon atoms, or, with R₃, forms an additional double bond so that the ethylenic bond becomes acetylenic, making it possible to go from an allyl radical to a propargyl radical;
- R₅ represents a hydrogen or a hydrocarbon-based radical such as an aryl radical or an alkyl, including aralkyl, radical, preferably of 1 or

2 carbon atoms;

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- R_5 and R_4 may be fractions of said "Ar" group above, such that R_5 and R_4 , and also the carbon which bears them, form a radical Ar;
- 5 it being possible for one of R_1 , R_2 , R_4 , R_3 and R_5 to be, in addition, chosen from specific trivalent, nitrile or acid functions, optionally and preferably in esterified form;
 - R' may be a hydrogen, a protective group, an aryl or an alkyl, including aralkyl, advantageously chiral;
 - R" is an allyl or homoallyl radical.
- 14. The use as claimed in claim 13, characterized in that just one of R' and R" is a homoallyl or allyl radical.